## Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

## Listing of Claims:

1. (Previously presented) A compound of Formula I,



or a pharmaceutically acceptable salt thereof, wherein,

Ar is selected from the following formulae



wherein Ar is substituted with -X and -Y-L-Z, in an ortho relationship to each other, and said Ar is optionally substituted with up to four R<sup>1</sup>;

each  $R^1$  is independently selected from -H, halogen, -CN, -NO<sub>2</sub>, -OR<sup>3</sup>, -N(R<sup>3</sup>)R<sup>3</sup>, -S(O)<sub>0-2</sub>R<sup>3</sup>, -SO<sub>2</sub>N(R<sup>3</sup>)R<sup>3</sup>, -CO<sub>2</sub>R<sup>3</sup>, -C(O)N(R<sup>3</sup>)R<sup>3</sup>, -N(R<sup>3</sup>)SO<sub>2</sub>R<sup>3</sup>, -N(R<sup>3</sup>)C(O)R<sup>3</sup>, -N(R<sup>3</sup>)CO<sub>2</sub>R<sup>3</sup>, -C(O)R<sup>3</sup>, -OC(O)R<sup>3</sup>, optionally substituted lower alkyl, optionally substituted aryl, optionally substituted lower arylalkyl, optionally substituted heterocyclyl, and optionally substituted lower heterocyclylalkyl;

X is selected from the following formulae.

wherein R4a is -C(O)N(R3)R3:

n = 1 or 2:

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p = 0 or 1;

a is 1 to 3:

M is  $-OR^3$  or  $-N(R^3)R^4$ ;

each  $R^2$  is independently selected from -H, haloalkyl, -C<sub>L6</sub>alkyl-N( $R^3$ ) $R^3$ , -C<sub>L6</sub>alkyl-OR $^3$ , -C<sub>L6</sub>alkyl-CO<sub>2</sub> $R^3$ , and -C<sub>L6</sub>alkyl-C(O)N( $R^3$ ) $R^3$ ;

cach R<sup>3</sup> is independently selected from -H, optionally substituted lower alkyl, optionally substituted aryl, optionally substituted lower arylalkyl, optionally substituted heterocyclyl, and optionally substituted lower heterocyclylalkyl; or

two of R<sup>3</sup>, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl ring, said optionally substituted five- to seven-membered heterocyclyl ring optionally containing at least one additional heteroatom selected from N, O, S, and P;

each  $R^4$  is independently selected from  $R^3$ ,  $-SO_2R^3$ ,  $-SO_2N(R^3)R^3$ ,  $-CO_2R^3$ ,  $-C(O)N(R^3)R^3$ , and  $-C(O)R^3$ ;

Y-L-Z is selected from the following formulae,

wherein g is zero to two; T is selected from absent, -N(R<sup>3</sup>)-, -S- and -O-; and each methylene between Y and T is optionally substituted; provided that when both Y and T are heteroatoms then g must be two;

Y is -O- or optionally substituted -CH2-;

 $R^5$  is selected from -H, halogen, -CN, -NO<sub>2</sub>, -OR<sup>3</sup>, -N( $R^3$ )R<sup>4</sup>, -S(O)<sub>0.2</sub>R<sup>3</sup>, -SO<sub>2</sub>N( $R^3$ )R<sup>3</sup>, -CO<sub>2</sub>R<sup>3</sup>, -C(O)N( $R^3$ )R<sup>3</sup>, -N( $R^3$ )SO<sub>2</sub>R<sup>3</sup>, -N( $R^3$ )C(O)R<sup>3</sup>, -N( $R^3$ )CO<sub>2</sub>R<sup>3</sup>, -C(O)R<sup>3</sup>, optionally substituted lower alkyl, optionally substituted aryl, optionally

substituted lower arylalkyl, optionally substituted heterocyclyl, and optionally substituted lower heterocyclylalkyl; and

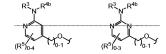
optionally two of R<sup>5</sup>, together with the atoms to which they are attached, form a second ring system fused with said five- to seven-membered ring system, said second ring system substituted with zero to four of R<sup>5</sup>.

- (Canceled)
- (Canceled)
- 4. (Canceled)
- 5. (Canceled)
- (Canceled)
- 7. (Canceled)
- 8. (Canceled)

(Canceled)

9.

- (Canceled)
- 11. (Canceled)
- 12. (Previously presented) The compound according to claim 1, wherein g is one or two.
- (Original) The compound according to claim 12, wherein each R<sup>5</sup> is independently selected from -H, halogen, -CN, -NH<sub>2</sub>, -NO<sub>2</sub>, -OR<sup>3</sup>, -N(R<sup>3</sup>)R<sup>4</sup>, -S(O)<sub>0-2</sub>R<sup>3</sup>, -SO<sub>2</sub>N(R<sup>3</sup>)R<sup>3</sup>, -CO<sub>2</sub>R<sup>3</sup>, -C(O)N(R<sup>3</sup>)R<sup>3</sup>, -N(R<sup>3</sup>)SO<sub>2</sub>R<sup>3</sup>, -N(R<sup>3</sup>)C(O)R<sup>3</sup>, -N(R<sup>3</sup>)CO<sub>2</sub>R<sup>3</sup>, -C(O)R<sup>3</sup>, and optionally substituted lower alkyl.
- (Currently amended) The compound according to claim 13, wherein -Y-L-Z is selected from the following formulae.



$$R^{3}_{N}$$
,  $R^{4b}$   $R^{2}_{N}$ ,  $R^{4b}$ 

wherein R<sup>4b</sup> is selected from R<sup>3</sup>, H, CO<sub>2</sub>R<sup>3</sup>, C(O)N(R<sup>3</sup>)R<sup>3</sup>, and C(O)R<sup>3</sup>.

15. (Previously presented) The compound according to claim 13, having formula III,

$$\mathbb{R}^{34}$$
 $\mathbb{R}^{5}$ 
 $\mathbb{R}^{5}$ 
 $\mathbb{R}^{5}$ 
 $\mathbb{R}^{5}$ 
 $\mathbb{R}^{36}$ 
 $\mathbb{R}^{46}$ 

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wherein J is N, and B is =N- or  $=C(R^5)-$ .

R<sup>3a</sup> is selected from optionally substituted aryl, optionally substituted lower arylalkyl, optionally substituted heterocyclyl, and optionally substituted lower heterocyclylalkyl;

R3b is H. and

$$R^{4b}$$
 is  $R^3$ , -H, -CO<sub>2</sub> $R^3$ , -C(O)N( $R^3$ ) $R^4$ , or -C(O) $R^3$ .

- 16. (Canceled)
- (Previously presented) The compound according to claim 15, wherein R<sup>3a</sup> is selected from optionally substituted aryl and optionally substituted heteroaryl.
- (Original) The compound according to claim 17, wherein R<sup>3a</sup> is optionally substituted phenyl.
- (Original) The compound according to claim 18, wherein said optionally substituted phenyl is substituted with at least one of halogen, -CN, -CF<sub>3</sub>, -NH<sub>2</sub>, -NO<sub>2</sub>, -OR<sup>3</sup>, -N(R<sup>3</sup>)R<sup>3</sup>, -S(O)<sub>0.2</sub>R<sup>3</sup>, -SO<sub>2</sub>N(R<sup>3</sup>)R<sup>3</sup>, -CO<sub>2</sub>R<sup>3</sup>, -CO<sub>2</sub>R<sup>3</sup>, -CO<sub>3</sub>R<sup>3</sup>, -CO<sub>3</sub>R<sup>3</sup>, -N(R<sup>3</sup>)SO<sub>2</sub>R<sup>3</sup>,

-N( $R^3$ )C(O) $R^3$ , -N( $R^3$ )CO $_2R^3$ , -C(O) $R^3$ , optionally substituted lower alkyl, and optionally substituted aryl.

- (Original) The compound according to claim 19, wherein said optionally substituted
  phenyl group is substituted with at least one trifluoromethyl group.
- (Original) The compound according to claim 20, wherein said optionally substituted
  phenyl group is substituted with at least two trifluoromethyl groups
- (Original) The compound according to claim 19, wherein said optionally substituted
  phenyl group is substituted with at least one lower alkyl group.
- 23. (Original) The compound according to claim 19, wherein R<sup>3b</sup> is -H.
- (Original) The compound according to claim 23, wherein R<sup>4b</sup> is selected from R<sup>3</sup>, -H,
  -CO<sub>2</sub>R<sup>3</sup>, -C(O)N(R<sup>3</sup>)R<sup>4</sup>, and -C(O)R<sup>3</sup>.
- (Canceled)
- (Canceled)
- (Original) The compound according to claim 24, wherein Ar is according to the formula below.

$$\overline{}$$

 (Original) The compound according to claim 24, wherein Ar is according to the formula below.

- (Canceled)
- 30. (Previously presented) A compound of Formula IV,

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$$(R^{6})_{0.5} \qquad (R^{5})_{0.3} \qquad H$$

$$(R^{6})_{0.5} \qquad (R^{2})_{0.4} \qquad B$$

or a pharmaceutically acceptable salt thereof, wherein,

Ar is selected from the following formulae:

each  $R^1$  is independently selected from -H, halogen, -CN, -NO<sub>2</sub>, -OR<sup>3</sup>, -N(R<sup>3</sup>)R<sup>3</sup>, -S(O)<sub>0-2</sub>R<sup>3</sup>, -SO<sub>2</sub>N(R<sup>3</sup>)R<sup>3</sup>, -CO<sub>2</sub>R<sup>3</sup>, -C(O)N(R<sup>3</sup>)R<sup>3</sup>, -N(R<sup>3</sup>)SO<sub>2</sub>R<sup>3</sup>, -N(R<sup>3</sup>)C(O)R<sup>3</sup>, -N(R<sup>3</sup>)CO<sub>2</sub>R<sup>3</sup>, -C(O)R<sup>3</sup>, optionally substituted lower alkyl, optionally substituted aryl, optionally substituted lower arylalkyl, optionally substituted heterocyclyl, and optionally substituted lower heterocyclylalkyl:

optionally two of  $\mathbb{R}^1$ , together with the atoms to which they are attached, form a first ring system fused with Ar, said first ring system substituted with zero to three additional of  $\mathbb{R}^1$ :

each  $R^2$  is independently selected from -H, halogen, oxo, -CN, -NH<sub>2</sub>, -NO<sub>2</sub>, -OR<sup>3</sup>, -N(R<sup>3</sup>)R<sup>3</sup>, -N(R<sup>3</sup>)R<sup>5</sup>, -S(O) $_{0}$ 2R<sup>3</sup>, -SO<sub>2</sub>N(R<sup>3</sup>)R<sup>3</sup>, -CO<sub>2</sub>R<sup>3</sup>, -C(O)N(R<sup>3</sup>)R<sup>3</sup>, -N(R<sup>3</sup>)SO<sub>2</sub>R<sup>3</sup>, -N(R<sup>3</sup>)C(O)R<sup>3</sup>, -N(R<sup>3</sup>)C(O)R<sup>3</sup>, -N(R<sup>3</sup>)C(O)R<sup>3</sup>, optionally substituted lower alkyl, optionally substituted aryl, optionally substituted lower arylalkyl, optionally substituted heterocyclyl, and optionally substituted lower heterocyclylalkyl;

two of R<sup>2</sup>, together with the atoms to which they are attached, can form an optionally substituted three- to seven-membered ring system;

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each R<sup>3</sup> is independently selected from -H, optionally substituted lower alkyl, optionally substituted arvl, optionally substituted lower arylalkyl, optionally substituted heterocyclyl, and optionally substituted lower heterocyclylalkyl; or

two of R3. when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl ring, said optionally substituted five- to seven-membered heterocyclyl ring optionally containing at least one additional heteroatom selected from N. O. S. and P:

each R4 is independently selected from R3, -SO2R3, -SO2N(R3)R3, -CO2R3, -C(O)N(R3)R3, and -C(O)R3;

Y is selected from optionally substituted -CH<sub>2</sub>-, -O-, -S-, and -N(R<sup>3</sup>)-;

L is selected from optionally substituted -CH<sub>2</sub>-, -O-, -S-, -N(R<sup>3</sup>)- and absent;

provided that Y and L are not both heteroatoms:

B is =N- or =C(H)-:

at each instance. R5 and R6 are independently selected from -H, halogen, -CN, -NO2,  $-OR^3$ ,  $-N(R^3)R^4$ ,  $-S(O)_{0.7}R^3$ ,  $-SO_2N(R^3)R^3$ ,  $-CO_2R^3$ ,  $-C(O)N(R^3)R$ ,  $-N(R^3)SO_2R^3$ . -N(R<sup>3</sup>)C(O)R<sup>3</sup>, -N(R<sup>3</sup>)CO<sub>2</sub>R<sup>3</sup>, -C(O)R<sup>3</sup>, optionally substituted lower alkyl, optionally substituted aryl, optionally substituted lower arylalkyl, optionally substituted heterocyclyl, and optionally substituted lower heterocyclylalkyl; and

optionally two of R5, together with the atoms to which they are attached, form a ring system fused with the ring containing B according to formula IV, said ring system substituted with zero to two additional of R5.

- (Original) The compound according to claim 30, wherein Y is -O- and L is optionally 31. substituted -CH2-.
- (Original) The compound according to claim 31, wherein at least one of R<sup>6</sup> is optionally 32 substituted lower alkyl.
- 33. (Original) The compound according to claim 32, wherein said at least one optionally substituted lower alkyl is meta- to the piperazine urea function as depicted in formula IV.

- (Original) The compound according to claim 33, wherein R<sup>4a</sup> is selected from R<sup>3</sup>, -H, -CO<sub>2</sub>R<sup>3</sup>, -C(O)N(R<sup>3</sup>)R<sup>4</sup>, and -C(O)R<sup>3</sup>.
- (Original) The compound according to claim 34, wherein R<sup>4a</sup> is selected from -H, -CO<sub>2</sub>R<sup>3</sup>, -C(O)N(R<sup>3</sup>)R<sup>4</sup>, and -C(O)R<sup>3</sup>.
- 36. (Original) The compound according to claim 35, wherein -Y-L- is -OCH2-.
- 37. (Canceled)
- (Canceled)
- (Original) The compound according to claim 36, wherein Ar is according to the formula below.

 (Original) The compound according to claim 36, wherein Ar is according to the formula below.

$$\langle \rangle$$

- (Canceled)
- 42. (Previously presented) A compound selected from Table 4.

Table 4

97	N-[3,5-bis(trifluoromethyl)phenyl]-4- {3-[(pyridin-4-ylmethyl)oxy]pyridin-2- yl}piperazine-1-carboxamide	HN N N N N N N N N N N N N N N N N N N
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Table 4

103	N-(4-chlorophenyl)-4-{3-[(pyridin-4-ylmethyl)oxy]pyridin-2-yl}piperazine- 1-carboxamide	
105	N-(3-chlorophenyl)-4-{3-[(pyridin-4- ylmethyl)oxy]pyridin-2-yl}piperazine- 1-carboxamide	
142	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-[3,5- bis(trifluoromethyl)phenyl]piperazine- 1-carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
144	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-(3- ethylphenyl)piperazine-1-carboxamide	NH2
161	methyl [4-({[2-(4-{[(3- ethylphenyl)amino]carbonyl}piperazin- 1-yl)pyridin-3-yl]oxy}methyl)pyridin- 2-yl]carbamate	

Table 4

164	methyl [4-({[2-(4-{[(3-bromophenyl)amino]carbonyl}piperazi n-1-yl)pyridin-3- yl]oxy}methyl)pyridin-2-yl]carbamate	Br 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2
165	methyl {4-[({2-[4-({[3- (methyloxy)phenyl]amino}carbonyl)pi perazin-1-yl]pyridin-3- yl}oxy)methyl]pyridin-2-yl}carbamate	NH NN NH
166	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy pyridin-2-yl)-N-[3- (methyloxy)phenyl]piperazine-1- carboxamide	NH <sub>2</sub> N N N N N N N N N N N N N N N N N N N
167	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-[3-(1- methylethyl)phenyl]piperazine-1- carboxamide	NH <sub>2</sub> N N N NH
168	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-{3- [(trifluoromethyl)oxy]phenyl}piperazin e-1-carboxamide	NH <sub>2</sub> NH <sub>2</sub> NH <sub>3</sub>
169	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy]pyridin-2-yl)-N-[2- fluoro-5- (trifluoromethyl)phenyl]piperazine-1- carboxamide	NH <sub>2</sub>

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Table 4

170	N-(3-ethylphenyl)-4-[3-({[2-({[(3-ethylphenyl)amino)carbonyl}amino)pyridin-4-yl]methyl}oxy)pyridin-2-yl]piperazine-1-earboxamide	HN NH NH NH
171	N-(3-ethylphenyl)-4-(3-{[(2-{[(4- methylpiperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- l-carboxamide	
173	N-[3,5-bis(trifluoromethyl)phenyl]-4- (3-{[(2-{[(4-methylpiperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- l-carboxamide	N N N N N N N N N N N N N N N N N N N
174	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-[3- (trifluoromethyl)phenyl]piperazine-1- carboxamide	H <sub>2</sub> N N F <sub>3</sub> C
175	4-[3-({[2-(acetylamino)pyridin-4-yl]methyl}oxy)pyridin-2-yl]-N-(3-ethylphenyl)piperazine-1-carboxamide	NH NH NH

Table 4

176	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-(3- ethyl-4-fluorophenyl)piperazine-1- carboxamide	H <sub>2</sub> N N N NH
177	2-[4-(3-{[(2-aminopyrimidin-4-yl)methyl]oxy}pyridin-2-yl)piperazin- 1-yl]-N-[3,5- bis(trifluoromethyl)phenyl]acetamide	NNNNNNNNNNNNNNNNNNNNNNNNNNNNNNNNNNNNNN
178	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N- phenylpiperazine-1-carboxamide	H <sub>2</sub> N N N NH
179	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-(3- chloro-5-ethylphenyl)piperazine-1- carboxamide	H <sub>2</sub> N N N NH
180	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-(5- ethyl-2-fluorophenyl)piperazine-1- carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
181	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-(3- bromo-5-ethylphenyl)piperazine-1- carboxamide	H <sub>2</sub> N N N NH

Amendment

Table 4

182	2-(4-methylpiperazin-1-yl)ethyl [4- ({[2-(4-{[(3- ethylphenyl)amino)earbonyl}piperazin- 1-yl)pyridin-3-yl]oxy}methyl)pyridin- 2-yl]carbamate	
183	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy   pyridin-2-yl)-N-(3- chlorophenyl)piperazine-1- carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
184	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy j pyridin-2-yl)-N-(3- bromophenyl)piperazine-1- carboxamide	H <sub>2</sub> N N N NH
185	N-[4-({[2-(4-acetylpiperazin-1- yl)pyridin-3-yl]oxy}methyl)pyridin-2- yl]-2-(4-methylpiperazin-1- yl)acetamide	
186	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy]pyridin-2-yl)-N-(3- fluorophenyl)piperazine-1- carboxamide	H <sub>2</sub> N N N NH

Table 4

187	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy] pyridin-2-yl)-N-(4- fluorophenyl)piperazine-1- carboxamide	H <sub>2</sub> N N NH
188	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-(2- fluorophenyl)piperazine-1- carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
189	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-(3,5- diethylphenyl)piperazine-1- carboxamide	H <sub>2</sub> N N N NH
190	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}-5-bromopyridin-2-yl)- N-(3-ethylphenyl)piperazine-1- carboxamide	Br N N NH
191	N-methyl-4-(3-{[(2-{[(4- methylpiperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- 1-carboxamide	N N N N N N N N N N N N N N N N N N N
192	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-[2- chloro-5- (trifluoromethyl)phenyl]piperazine-1- carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
193	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy]pyridin-2-yl)-N-(5- chloro-2-fluorophenyl)piperazine-1- carboxamide	H <sub>2</sub> N N CI F

Table 4

194	4-(3-{[(2-amino-5-bromopyrimidin-4-yl)methyl]oxy}-5-bromopyridin-2-yl)-N-(3-ethylphenyl)piperazine-1-carboxamide	Br N N NH
195	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-[2- fluoro-3- (trifluoromethyl)phenyl]piperazine-1- carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
196	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy} pyridin-2-yl)-N-[3- fluoro-5- (trifluoromethyl)phenyl]piperazine-1- carboxamide	H <sub>2</sub> N N F CF <sub>3</sub>
197	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy} pyridin-2-yl)-N-(3,5- dichlorophenyl)piperazine-1- carboxamide	H <sub>2</sub> N N CI CI
198	N-(3-chloro-5-ethylphenyl)-4-(3-{[(2- {[(4-methylpiperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- 1-carboxamide	N N N N N N N N N N N N N N N N N N N

Table 4

199	N-(5-ethyl-2-fluorophenyl)-4-(3-{[(2- {[(4-methylpiperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- 1-carboxamide	N N N N N N N N N N N N N N N N N N N
200	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-[3- ethyl-5- (trifluoromethyl)phenyl]piperazine-1- carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
204	4-(3-{[(2-aminopyrimidin-4-yl)methyl]oxy}pyridin-2-yl)-N-methylpiperazine-1-carboxamide	H <sub>2</sub> N N NH
205	4-(3-{[(2-aminopyrimidin-4-yl)methyl]oxy}pyridin-2-yl)-N-ethylpiperazine-1-carboxamide	H <sub>2</sub> N N NH
206	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N- cyclohexylpiperazine-1-carboxamide	H <sub>2</sub> N N N NH
207	4-({[2-(4-acetylpiperazin-1-yl)pyridin-3-yl]oxy}methyl)pyrimidin-2-amine	O N N N N N N N N N N N N N N N N N N N
208	4-({[2-(4-propanoylpiperazin-1-yl)pyridin-3-yl]oxy}methyl)pyrimidin- 2-amine	N N N N N N N N N N N N N N N N N N N

Table 4

209	N-(3-cyclopropylphenyl)-4-(3-{[[2- {[[(4-methylpiperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- 1-carboxamide	N N N N N N N N N N N N N N N N N N N
210	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-(3- cyclopropylphenyl)piperazine-1- carboxamide	H <sub>2</sub> N N NH
211	N-[2-fluoro-5- (trifluoromethyl)phenyl]-4-(3-{[(2- {[(4-methyl)piperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- 1-carboxamide	N N N N N N N N N N N N N N N N N N N
212	N-[3-fluoro-5- (trifluoromethyl)phenyl]-4-(3-{[(2- {[(4-methyl)piperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- 1-carboxamide	N N N N N N N N N N N N N N N N N N N
213	N-(3,5-dichlorophenyl)-4-(3-{[(2-{[(4- methylpiperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- 1-carboxamide	N N N N N N N N N N N N N N N N N N N

Amendment

Table 4

214	4-(3-{[(2-{[(4-methylpiperazin-1-yl)acetyl]amino}pyridin-4-yl)methyl[oxy]pyridin-2-yl)-N-[3-(trifluoromethyl)piperazine-1-carboxamide	N N N N N N N N N N N N N N N N N N N
216	4-(3-{[1-(2-aminopyrimidin-4- yl)ethyl]oxy pyridin-2-yl)-N-[3,5- bis(trifluoromethyl)phenyl]piperazine- 1-carboxamide	$H_2N$ $N$ $F_3C$ $CF_3$
219	4-[({2-[4-(3,4-dihydroquinolin-1(2H)- ylcarbonyl)piperazin-1-yl]pyridin-3- yl}oxy)methyl]pyrimidin-2-amine	N N N N N N N N N N N N N N N N N N N
220	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-(2- methylpropyl)piperazine-1- carboxamide	H <sub>2</sub> N NH
226	N-(3,5-diethylphenyl)-4-(3-{[(2-{[(4-methylpiperazin-1-yl)acetyl]amino}pyridin-4-yl)oxy]methyl}pyridin-2-yl)piperazine-1-carboxamide	NH NH
227	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}-6-methylpyridin-2-yl)- N-(3-ethylphenyl)piperazine-1- carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N

Table 4

228	4-(3-{[(2-aminopyrimidin-4-yl)methyl]oxy}-6-methylpyridin-2-yl)-N-[3,5-bis(trifluoromethyl)phenyl]piperazine-1-carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
233	N-[3-chloro-5- (trifluoromethyl)phenyl]-4-(3-{[(2- {[(4-methyl)piperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- 1-carboxamide	N N N N N N N N N N N N N N N N N N N
235	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-[3- chloro-5- (trifluoromethyl)phenyl]piperazine-1- carboxamide	H <sub>2</sub> N N N NH NH CI
237	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}-6-chloropyridin-2-yl)- N-(3-ethylphenyl)piperazine-1- carboxamide	CI N N NH
243	N-[3-chloro-2-fluoro-5- (trifluoromethyl)phenyl]-4-(3-{[(2- {[(4-methyl)piperazin-1- yl)acetyl]amino}pyridin-4- yl)methyl]oxy}pyridin-2-yl)piperazine- 1-carboxamide	NH P <sub>3</sub> C F

Table 4

244	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}-6-chloropyridin-2-yl)- N-[3,5- bis(trifluoromethyl)phenyl]piperazine- 1-carboxamide	$H_2N$ $N$ $F_3C$ $CF_3$
245	4-(3-{[1-(2-aminopyrimidin-4-yl)ethyl]oxy}pyridin-2-yl)-N-(3-ethylphenyl)piperazine-1-carboxamide	H <sub>2</sub> N N N NH
246	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}-6-chloropyridin-2-yl)- N-(5-ethyl-2-fluorophenyl)piperazine- 1-carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
247	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-(3- ethyl-5-fluorophenyl)piperazine-1- earboxamide	H <sub>2</sub> N N N NH
249	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy}pyridin-2-yl)-N-[3- chloro-2-fluoro-5- (trifluoromethyl)phenyl]piperazine-1- carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N
250	4-(3-{[(2-aminopyrimidin-4- yl)methyl]oxy} pyridin-2-yl)-N-[3,5- bis(trifluoromethyl)phenyl]-N- methylpiperazine-1-carboxamide	H <sub>2</sub> N N N N N N N N N N N N N N N N N N N

 (Previously presented) A pharmaceutical composition comprising the compound according to claim 1 and a pharmaceutically acceptable carrier.

- 44. (Canceled)
- 45. (Canceled)
- 46. (Canceled)
- 47. (Canceled)
- 48. (Canceled)
- 49. (Canceled)
- 50. (Canceled)
- 51. (Canceled)
- 52. (Canceled)
- 53. (Canceled)
- 54. (Canceled)
- 55. (Canceled)
- 56. (Canceled)
- 57. (Canceled)